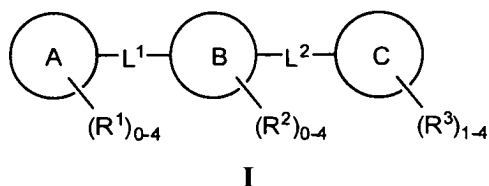


Amendments to the Claims:

This listing of claims will replace all prior versions, and listings, of claims in the application:

Listing of Claims:

1. (currently amended) A compound for modulating c-Kit activity according to Formula I,



or a pharmaceutically acceptable salt, thereof, wherein,

ring A is:

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wherein each Y is independently either =C(H)- or =N-; and Z is selected from -O-, -S-, and -N(R⁷)-;

each R^1 is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R^1 , together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{10} ;

L^1 is a single bond;

ring B is a five- to ten-membered aryl or a five- to ten-membered heterocyclyl;

each R^2 is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two adjacent of R^2 , together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{15} ;

L^2 is selected from -N(H)N(H)C(=O)N(H)-, -CH₂N(H)C(=O)N(H)-, -CH₂OC(=O)N(H)-, and -XCH₂C(=O)N(H)-; wherein X is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-; and any C-H of L^2 is optionally C-R²⁰;

ring C is phenyl or pyridyl;

each R^3 is independently selected from halogen, trihalomethyl, -CN,

-NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl

C₁₋₆alkyl; provided R³ is not a cyclic sulfonamide attached to ring C via the nitrogen of said cyclic sulfonamide;

two adjacent of R³, together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R²⁵;

R⁴ is selected from -H, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

two of R⁴, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R⁵ is selected from -H, -CN, -NO₂, -OR⁴, -S(O)₀₋₂R⁴, -CO₂R⁴, optionally substituted C₁₋₆alkyl, optionally substituted C₁₋₆alkenyl, and optionally substituted C₁₋₆alkynyl;

R⁷ is selected from -H, optionally substituted C₁₋₆alkyl, -SO₂N(R⁴)R⁴, -CO₂R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl; and

each of R¹⁰, each of R¹⁵, each of R²⁰, and each of R²⁵ is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴, -N(R⁴)R⁴, -S(O)₀₋₂R⁴, -SO₂N(R⁴)R⁴, -CO₂R⁴, -C(=O)N(R⁴)R⁴, -C(=NR⁵)N(R⁴)R⁴, -C(=NR⁵)R⁴, -N(R⁴)SO₂R⁴, -N(R⁴)C(O)R⁴, -NCO₂R⁴, -C(=O)R⁴, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

provided:

1) when both ring B and ring C are phenyl:

a) and the compound comprises ring B-CH₂N(H)C(=O)N(H)-ring C, then L¹ must be a single bond; R³ can not comprise a group of the formula

-O(CH₂)₂₋₄-N-piperazine that is *ortho*- to L²; and ring A cannot be a 5-methyl-[1,2,4]-oxadiazol-3-yl radical, a 4H-[1,2,4]-oxadiazol-5-one-3-yl radical, nor a 4'-[2,2';6',2'']terpyridinyl radical;

b) and L¹ is single bond, then L² cannot comprise -N(H)C(=O)C(=O)N(H)- nor -N(H)C(=Q)C(H)CNC(=O)- (where Q is S or O);

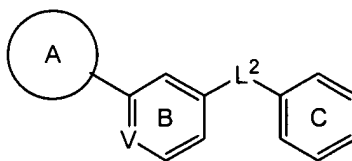
c) and L¹ is other than single bond, then A cannot be quinolin-2-yl-L¹, quinolin-3-yl-L¹, or quinolin-4-yl-L¹;

2) when ring A is a fused aryl system, then L¹ must be a single bond;

3) when ring B is phenyl, ring C is a C₆₋₁₆carbocyclic, L¹ is a single bond, and the compound comprises -ring B-OCH₂C(=O)N(H)- then ring A cannot be a 2,5-dimethyl-1H-pyrrole-1-yl radical;

4) ring A cannot be a pyrimidin-2-yl radical when L¹ is -N(H)- and ring B is phenyl;

5) when the compound comprises the formula,



where V is =C(H)- or =N-, and there is a nitrogen of L² bound directly to ring B, then A can not comprise a [1,2,4]-oxadiazol-3-yl radical; and

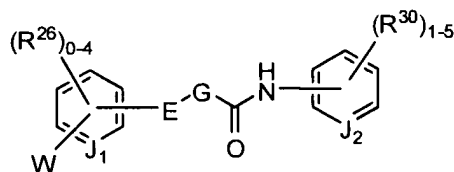
6) the compound is not one of: N-naphthalen-1-yl-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(phenyloxy)phenyl]-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,4-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,6-dimethylphenyl)-2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-

trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-(4-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-(2,6-diethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-[2-(ethyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-[4-(dimethylamino)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-(4-chloro-3-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 yl)phenyl]oxy}acetamide, N-(4-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 (trifluoromethyl)phenyl]acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoromethyl)phenyl]acetamide,
 (trifluoromethyl)phenyl]acetamide, methyl 4-[[{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate,
 ethyl 4-[[{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoate, 3-[[{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino]benzoic acid,
 N-[3-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 {{3-(4H-1,2,4-triazol-4-yl)phenyl]oxy}acetamide, N-(4-chlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide,
 N-(4-aminophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, and N-(4-acetylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide.

2-10. (Cancelled)

11. (original) The compound according to claim 10, wherein there exists at least one of R³ that is halogen.

12. (original) The compound according to claim 10, wherein there exists at least one of R^3 that is trihalomethyl.
13. (original) The compound according to claim 10, wherein there exists at least one of R^3 that is trifluoromethyl.
14. (original) The compound according to claim 13, wherein ring C is a phenyl comprising a trifluoromethyl radical *meta*- to L^2 .
15. (previously presented) The compound according to claim 10, wherein each of R^3 is independently selected from halogen, trihalomethyl, $-OR^4$, $-C(=O)R^4$, and optionally substituted C_{1-6} alkyl.
16. (currently amended) A compound for modulating c-Kit activity according to Formula II,



II

or a pharmaceutically acceptable salt, thereof, wherein,

W is selected from the following:

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each of R^{27} independently selected from halogen, trihalomethyl, -CN, -NO₂, -OR⁵⁵, -S(O)₀₋₂R⁵⁵, -SO₂N(R⁵⁵)R⁵⁵, -C(=O)N(R⁵⁵)R⁵⁵, -C(=NR⁵⁰)N(R⁵⁵)R⁵⁵, -C(=NR⁵⁰)R⁵⁵, -N(R⁵⁵)SO₂R⁵⁵, -N(R⁵⁵)C(O)R⁵⁵, -NCO₂R⁵⁵, -C(=O)R⁵⁵, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

each Y is independently either =C(H)- or =N-;

Z is selected from -O-, -S(O)₀₋₂-, and -N(R⁷)-

E and G are each independently selected from -O-, -S(O)₀₋₂-, -C(R³¹)R³²-, and -N(R³³)-

J₁ and J₂ are each independently =C(H)- or =N-;

R²⁶ is independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -CO₂R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰, -N(R⁴⁰)SO₂R⁴⁰, -N(R⁴⁰)C(O)R⁴⁰, -NCO₂R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl;

R³⁰ is independently selected from halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -S(O)₀₋₂R⁴⁰, -SO₂N(R⁴⁰)R⁴⁰, -C(=O)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)N(R⁴⁰)R⁴⁰, -C(=NR⁵⁰)R⁴⁰,

$-N(R^{40})SO_2R^{40}$, $-N(R^{40})C(O)R^{40}$, $-NCO_2R^{40}$, $-C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl; or

two adjacent of R^{26} or two adjacent of R^{30} , together with the annular atoms to which they are attached, can form a five- to six-membered ring containing up to two heteroatoms and optionally substituted with up to three of R^{35} ;

R^{31} and R^{32} are each independently selected from -H, halogen, trihalomethyl, -CN, -NO₂, -OR⁴⁰, $-N(R^{40})R^{40}$, $-S(O)_{0-2}R^{40}$, $-SO_2N(R^{40})R^{40}$, $-CO_2R^{40}$, $-C(=O)N(R^{40})R^{40}$, $-C(=NR^{50})N(R^{40})R^{40}$, $-C(=NR^{50})R^{40}$, $-N(R^{40})SO_2R^{40}$, $-N(R^{40})C(O)R^{40}$, $-NCO_2R^{40}$, $-C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl;

R^{33} is selected from -H, optionally substituted lower alkyl, $-SO_2N(R^{40})R^{40}$, $-CO_2R^{40}$, $-C(=O)N(R^{40})R^{40}$, $-C(=NR^{50})N(R^{40})R^{40}$, $-C(=NR^{50})R^{40}$, $-C(=O)R^{40}$, optionally substituted alkoxy, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl;

R^{40} is selected from -H, optionally substituted alkoxy, optionally substituted C_{1-6} alkyl, optionally substituted aryl, optionally substituted aryl C_{1-6} alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C_{1-6} alkyl;

two of R^{40} , when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P;

R^{50} is selected from -H, -CN, -NO₂, -OR⁴⁰, $-S(O)_{0-2}R^{40}$, $-CO_2R^{40}$, optionally substituted C_{1-6} alkyl, optionally substituted C_{1-6} alkenyl, and optionally substituted C_{1-6} alkynyl;

R⁵⁵ is selected from -H, optionally substituted C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl; and

two of R⁵⁵, when taken together with a common nitrogen to which they are attached, form an optionally substituted five- to seven-membered heterocyclyl, said optionally substituted five- to seven-membered heterocyclyl optionally containing at least one additional heteroatom selected from N, O, S, and P.

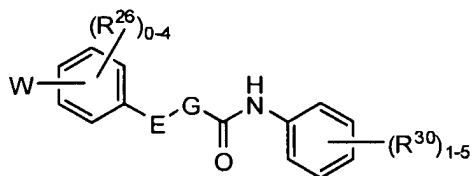
17. (original) The compound according to claim 16, wherein the annular carbons of ring B to which W and E are attached are not contiguous.

18. (previously presented) The compound according to claim 17, wherein R³⁰ is selected from halogen, trihalomethyl, -OR⁴⁰, -N(R⁴⁰)R⁴⁰, -C(=O)R⁴⁰, optionally substituted alkoxy, optionally substituted C₁₋₆alkyl, optionally substituted aryl, optionally substituted aryl C₁₋₆alkyl, optionally substituted heterocyclyl, and optionally substituted heterocyclyl C₁₋₆alkyl.

19. (original) The compound according to claim 18, wherein there exists at least one of R³⁰ that is trihalomethyl.

20. (original) The compound according to claim 18, wherein there exists at least one of R³⁰ that is trifluoromethyl.

21. (previously presented) The compound according to claim 18, according to formula III.



III

22. (currently amended) The compound according to claim 21, wherein W is selected from the

following:

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and R^{27} is defined as above.

23. (withdrawn) The compound according to claim 22, wherein E is selected from -O-, -S(O)₀₋₂-, and -NH-; and G is -CH₂-.

24. (withdrawn) The compound according to claim 22, wherein E is either -CH₂- or -NH-; and G is selected from -O-, -S-, and -NH-.

25. (Cancelled)

26. (original) The compound according to claim 25, wherein at least one of R^{30} is a trifluoromethyl radical *meta*- to -E-G-C(=O)N(H)-.

27. (currently amended) The compound according to claim 1, selected from Table 3:

Table 3

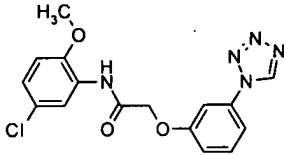
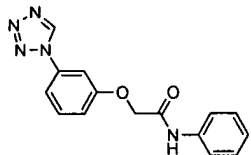
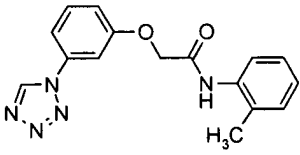
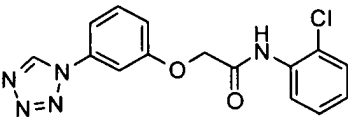
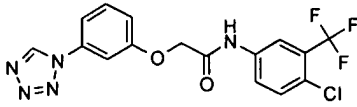
| Entry | Name | Structure |
|-------|---|--|
| 1 | N-[5-chloro-2-(methoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 2 | N-phenyl-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 3 | N-(2-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 4 | N-(2-chlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 5 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |

Table 3

| Entry | Name | Structure |
|-------|---|-----------|
| 6 | ethyl 2-[(3-(1H-tetrazol-1-yl)phenyl]oxy} acetyl)amino]-4,5,6,7-tetrahydro-1-benzothiophene-3-carboxylate | |
| 7 | N-(3-chloro-2-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |
| 8 | N-(3-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |
| 9 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(2H-tetrazol-5-yl)phenyl]oxy}acetamide | |
| 10 | N-(4-chloro-2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |

Table 3

| Entry | Name | Structure |
|-------|---|-----------|
| 11 | N-(4-bromo-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |
| 12 | N-(4-morpholin-4-ylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |
| 13 | N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |
| 14 | N-[4-bromo-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |
| 15 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |

Table 3

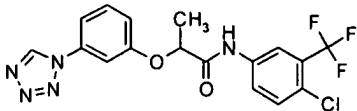
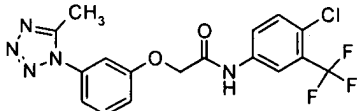
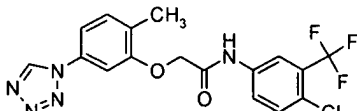
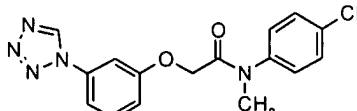
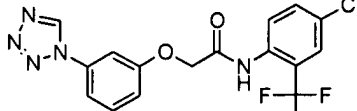
| Entry | Name | Structure |
|-------|--|--|
| 16 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}propanamide |  |
| 17 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(5-methyl-1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 18 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-methyl-5-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 19 | N-(4-chlorophenyl)-N-methyl-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 20 | N-[4-chloro-2-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |

Table 3

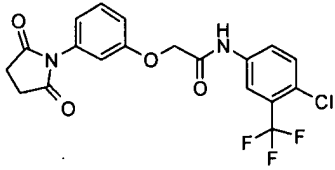
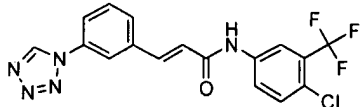
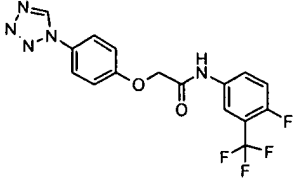
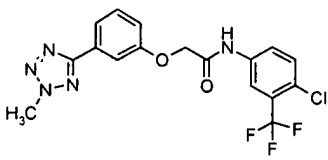
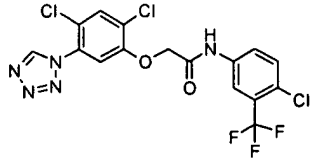
| Entry | Name | Structure |
|-------|--|--|
| 21 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(2,5-dioxopyrrolidin-1- yl)phenyl]oxy}acetamide |  |
| 22 | (2E)-N-[4-chloro-3- (trifluoromethyl)phenyl]-3-[3-(1H- tetrazol-1-yl)phenyl]prop-2-enamide |  |
| 23 | N-[4-fluoro-3-(trifluoromethyl)phenyl]-2- {[4-(1H-tetrazol-1- yl)phenyl]oxy}acetamide |  |
| 24 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(2-methyl-2H-tetrazol-5- yl)phenyl]oxy}acetamide |  |
| 25 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2,4-dichloro-5-(1H-tetrazol-1- yl)phenyl]oxy}acetamide |  |

Table 3

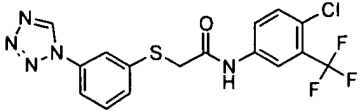
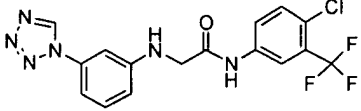
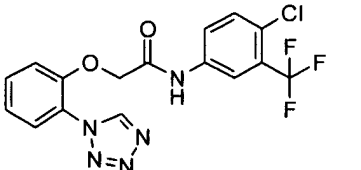
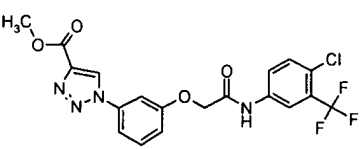
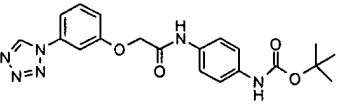
| Entry | Name | Structure |
|-------|--|--|
| 26 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(1H-tetrazol-1- yl)phenyl]thio}acetamide |  |
| 27 | N-[4-chloro-3-(trifluoromethyl)phenyl]- N-2-~[3-(1H-tetrazol-1- yl)phenyl]glycinamide |  |
| 28 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[2-(1H-tetrazol-1- yl)phenyl]oxy}acetamide |  |
| 29 | methyl 1-{3-[(2-{[4-chloro-3- (trifluoromethyl)phenyl]amino}-2- oxoethyl)oxy]phenyl}-1H-1,2,3-triazole-4- carboxylate |  |
| 30 | 1,1-dimethylethyl {4-[[[3-(1H-tetrazol-1- yl)phenyl]oxy}acetyl]amino]phenyl}carba mate |  |

Table 3

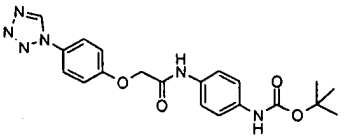
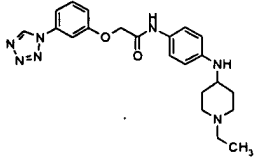
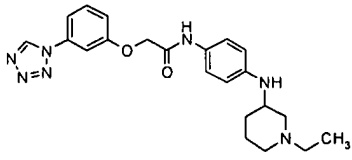
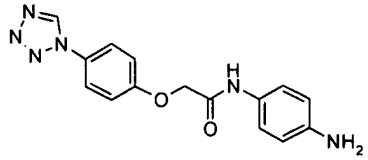
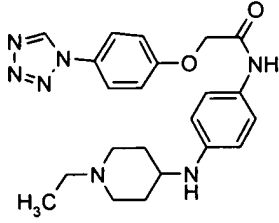
| Entry | Name | Structure |
|-------|---|---|
| 31 | 1,1-dimethylethyl {4-[[[4-(1H-tetrazol-1-yl)phenyl]oxy}acetyl]amino]phenyl} carbamate |  |
| 32 | N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 33 | N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 34 | N-(4-aminophenyl)-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 35 | N-{4-[(1-ethylpiperidin-4-yl)amino]phenyl}-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |

Table 3

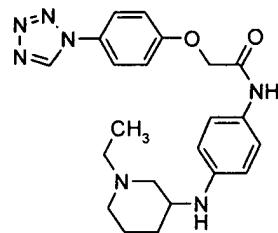
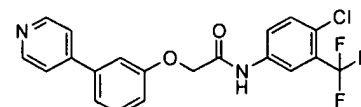
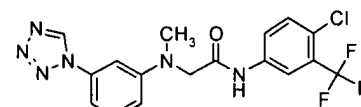
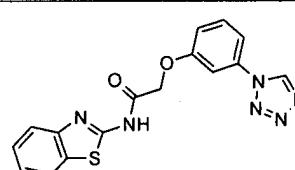
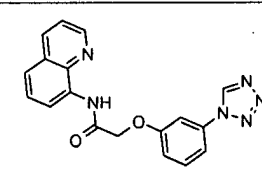
| Entry | Name | Structure |
|-------|---|--|
| 36 | N-{4-[(1-ethylpiperidin-3-yl)amino]phenyl}-2-{[4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 37 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-4-ylphenyl)oxy]acetamide |  |
| 38 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-methyl-N~2~-{3-(1H-tetrazol-1-yl)phenyl}glycinamide |  |
| 39 | N-1,3-benzothiazol-2-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 40 | N-quinolin-8-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |

Table 3

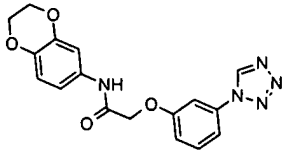
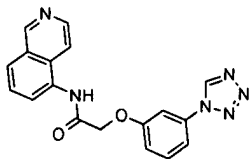
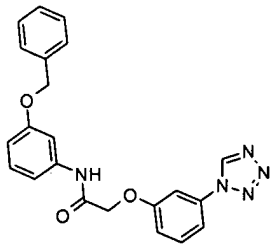
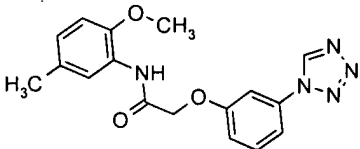
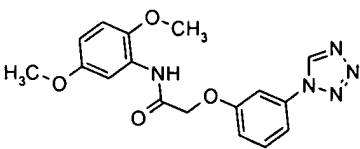
| Entry | Name | Structure |
|-------|---|--|
| 41 | N-(2,3-dihydro-1,4-benzodioxin-6-yl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 42 | N-isoquinolin-5-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 43 | N-{3-[(phenylmethyl)oxy]phenyl}-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 44 | N-[5-methyl-2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |
| 45 | N-[2,5-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide |  |

Table 3

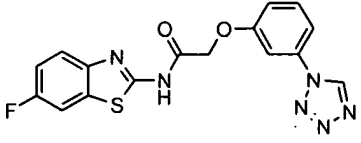
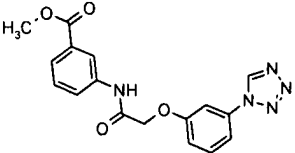
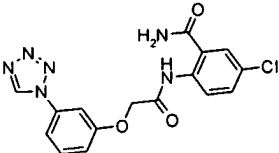
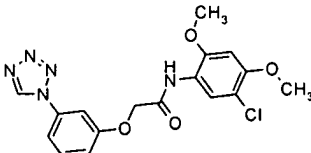
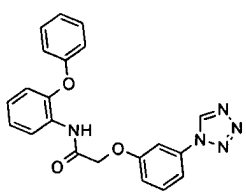
| Entry | Name | Structure |
|-------|--|--|
| 46 | N-(6-fluoro-1,3-benzothiazol-2-yl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 47 | methyl 3-(((3-(1H-tetrazol-1-yl)phenyl)oxy)acetyl)amino]benzoate |  |
| 48 | 5-chloro-2-(((3-(1H-tetrazol-1-yl)phenyl)oxy)acetyl)amino]benzamide |  |
| 49 | N-[5-chloro-2,4-bis(methoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 50 | N-[2-(phenyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |

Table 3

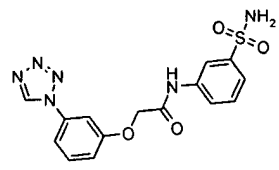
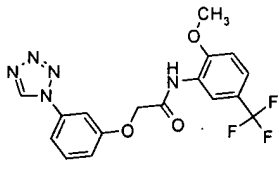
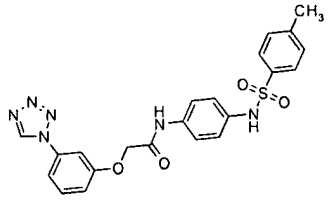
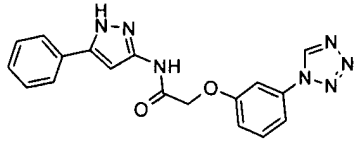
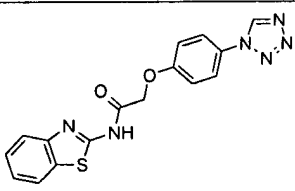
| Entry | Name | Structure |
|-------|--|--|
| 51 | N-[3-(aminosulfonyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 52 | N-[2-(methoxy)-5-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 53 | N-(4-{{(4-methylphenyl)sulfonyl}amino}phenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 54 | N-(5-phenyl-1H-pyrazol-3-yl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |
| 55 | N-1,3-benzothiazol-2-yl-2-{{4-(1H-tetrazol-1-yl)phenyl}oxy}acetamide |  |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 56 | N-quinolin-8-yl-2-{{4-(1H-tetrazol-1-yl)phenyl}oxy}acetamide | |
| 57 | 1,1-dimethylethyl 2-{3-[(2-{{4-chloro-3-(trifluoromethyl)phenyl}amino}-2-oxoethyl)oxy]phenyl}-1H-pyrrole-1-carboxylate | |
| 58 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-pyrrol-2-yl)phenyl}oxy}acetamide | |
| 59 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyrimidin-5-yl)phenyl]oxyacetamide | |
| 60 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-1,2,3-triazol-1-yl)phenyl}oxy}acetamide | |
| 61 | 4-chloro-N-(2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}ethyl)-3-(trifluoromethyl)aniline | |

Table 3

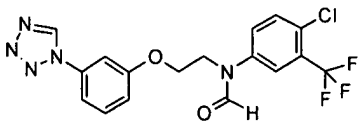
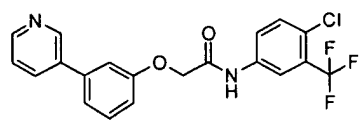
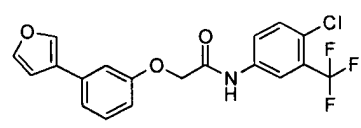
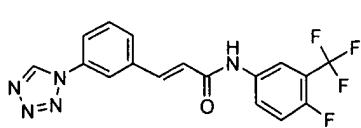
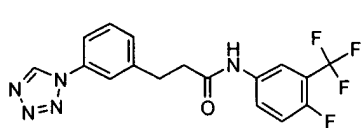
| Entry | Name | Structure |
|-------|--|--|
| 62 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N-(2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}ethyl)formamide |  |
| 63 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-pyridin-3-ylphenyl)oxy]acetamide |  |
| 64 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-furan-3-ylphenyl)oxy]acetamide |  |
| 65 | (2E)-N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]prop-2-enamide |  |
| 66 | N-[4-fluoro-3-(trifluoromethyl)phenyl]-3-[3-(1H-tetrazol-1-yl)phenyl]propanamide |  |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 67 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[6-(1H-tetrazol-1-yl)pyrimidin-4- yl]oxy}acetamide | |
| 68 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[3-(3,5-dimethylisoxazol-4- yl)phenyl]oxy}acetamide | |
| 69 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-quinolin-7-ylphenyl)oxy]acetamide | |
| 70 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [(3-furan-2-ylphenyl)oxy]acetamide | |
| 71 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- [3-(1H-tetrazol-1- yl)phenyl]hydrazinecarboxamide | |

Table 3

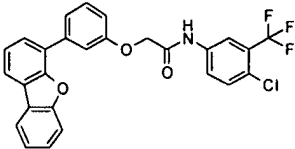
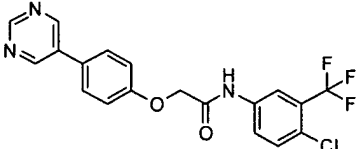
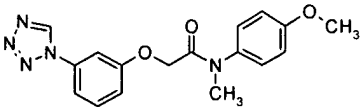
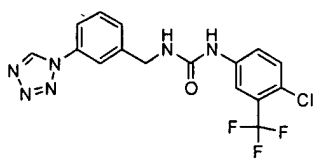
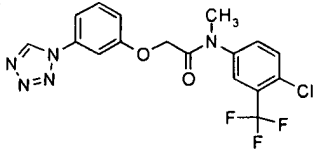
| Entry | Name | Structure |
|-------|---|--|
| 72 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(3-dibenzo[b,d]furan-4-ylphenyl)oxy]acetamide |  |
| 73 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-ylphenyl)oxy]acetamide |  |
| 74 | N-methyl-N-[4-(methoxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide |  |
| 75 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl} urea |  |
| 76 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N-methyl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide |  |

Table 3

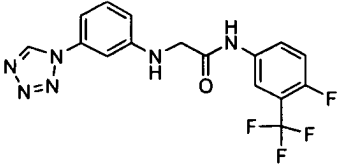
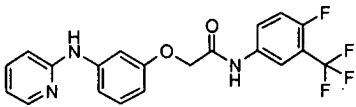
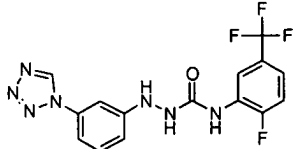
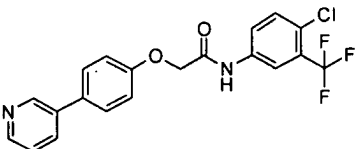
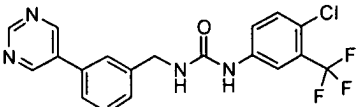
| Entry | Name | Structure |
|-------|---|--|
| 77 | N-[4-fluoro-3-(trifluoromethyl)phenyl]-N~2~-{3-(1H-tetrazol-1-yl)phenyl}glycinamide |  |
| 78 | N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-{{3-(pyridin-2-ylamino)phenyl}oxy}acetamide |  |
| 79 | N-[2-fluoro-5-(trifluoromethyl)phenyl]-2-[3-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide |  |
| 80 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide |  |
| 81 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea |  |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 82 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyrimidin-5-yl)phenyl)methyl]urea | |
| 83 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-yl)phenyl)methyl]urea | |
| 84 | [3-(1H-tetrazol-1-yl)phenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 85 | N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyrimidin-5-yl)phenyl]oxy]acetamide | |
| 86 | N-[2-[[4-chloro-3-(trifluoromethyl)phenyl]-N-[3-(1H-tetrazol-1-yl)phenyl]glycinamide | |

Table 3

| Entry | Name | Structure |
|-------|---|-----------|
| 87 | 2-{[4-chloro-3-(trifluoromethyl)phenyl]oxy}-N-[3-(1H-tetrazol-1-yl)phenyl]acetamide | |
| 88 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-methyl-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |
| 89 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[4-(1H-1,2,3-triazol-1-yl)phenyl]oxy}acetamide | |
| 90 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[3-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |
| 91 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{[2-fluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |

Table 3

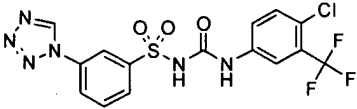
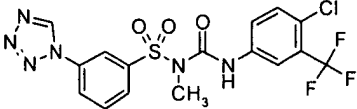
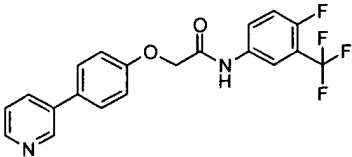
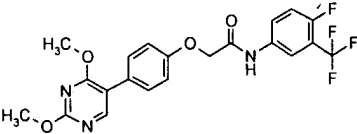
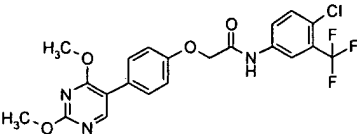
| Entry | Name | Structure |
|-------|---|--|
| 92 | N-([4-chloro-3-(trifluoromethyl)phenyl]amino)carbonyl-3-(1H-tetrazol-1-yl)benzenesulfonamide |  |
| 93 | N-([4-chloro-3-(trifluoromethyl)phenyl]amino)carbonyl-N-methyl-3-(1H-tetrazol-1-yl)benzenesulfonamide |  |
| 94 | N-[4-fluoro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-3-ylphenyl)oxy]acetamide |  |
| 95 | 2-([4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl]oxy)-N-[4-fluoro-3-(trifluoromethyl)phenyl]acetamide |  |
| 96 | 2-([4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl]oxy)-N-[4-chloro-3-(trifluoromethyl)phenyl]acetamide |  |

Table 3

| Entry | Name | Structure |
|-------|---|-----------|
| 97 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[(4-pyridin-4-ylphenyl)oxy]acetamide | |
| 98 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N-2-[3-(methoxy)-4-(1H-tetrazol-1-yl)phenyl]glycinamide | |
| 99 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N-2-[4-(methoxy)-3-(1H-tetrazol-1-yl)phenyl]glycinamide | |
| 100 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N-2-[4-(1H-tetrazol-1-yl)phenyl]glycinamide | |
| 101 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(2,3,5,6-tetrafluoro-4-pyrimidin-5-ylphenyl)hydrazinecarboxamide | |

Table 3

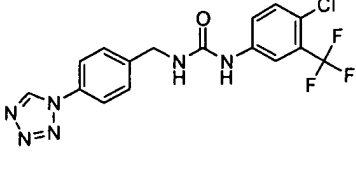
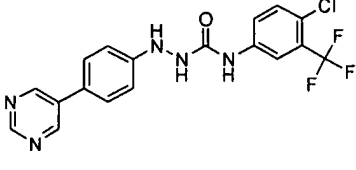
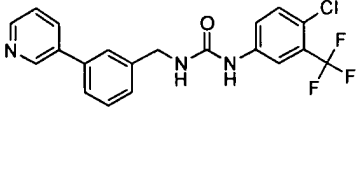
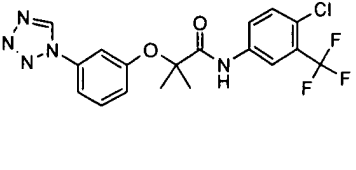
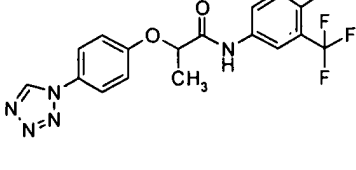
| Entry | Name | Structure |
|-------|---|--|
| 102 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(1H-tetrazol-1-yl)phenyl]methyl} urea |  |
| 103 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- (4-pyrimidin-5- ylphenyl)hydrazinecarboxamide |  |
| 104 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- [(3-pyridin-3-ylphenyl)methyl]urea |  |
| 105 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- methyl-2-{[3-(1H-tetrazol-1- yl)phenyl]oxy} propanamide |  |
| 106 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(1H-tetrazol-1- yl)phenyl]oxy} propanamide |  |

Table 3

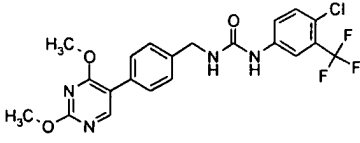
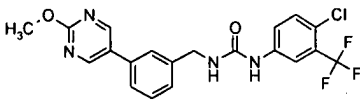
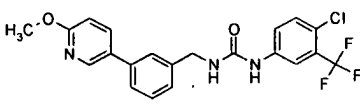
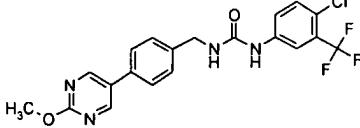
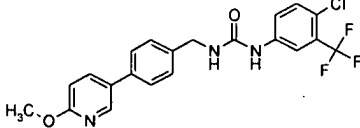
| Entry | Name | Structure |
|-------|--|--|
| 107 | N-({4-[2,4-bis(methyloxy)pyrimidin-5-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea |  |
| 108 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea |  |
| 109 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea |  |
| 110 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[2-(methyloxy)pyrimidin-5-yl]phenyl}methyl)urea |  |
| 111 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({4-[6-(methyloxy)pyridin-3-yl]phenyl}methyl)urea |  |

Table 3

| Entry | Name | Structure |
|-------|---|-----------|
| 112 | 1,1-dimethylethyl 2-{4-[(2-{[4-chloro-3-(trifluoromethyl)phenyl]amino}-2-oxoethyl)oxy]phenyl}-1H-indole-1-carboxylate | |
| 113 | N-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)-4-(1H-tetrazol-1-yl)benzenesulfonamide | |
| 114 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N~2~-[3-(2H-tetrazol-5-yl)phenyl]glycinamide | |
| 115 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[2,6-difluoro-4-(1H-tetrazol-1-yl)phenyl]oxy}acetamide | |
| 116 | (3-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |

Table 3

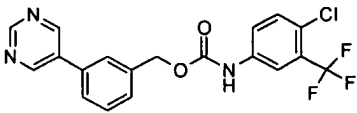
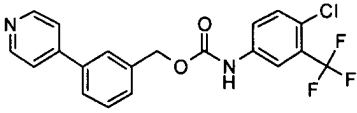
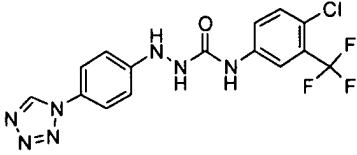
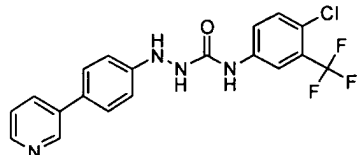
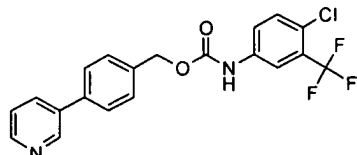
| Entry | Name | Structure |
|-------|---|--|
| 117 | (3-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |
| 118 | (3-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |
| 119 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-[4-(1H-tetrazol-1-yl)phenyl]hydrazinecarboxamide |  |
| 120 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-3-ylphenyl)hydrazinecarboxamide |  |
| 121 | (4-pyridin-3-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |

Table 3

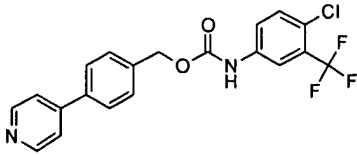
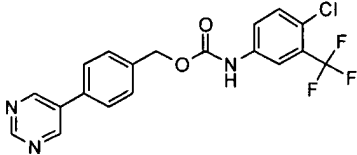
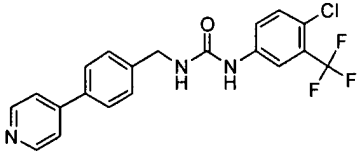
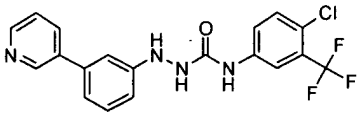
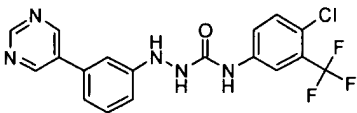
| Entry | Name | Structure |
|-------|---|--|
| 122 | (4-pyridin-4-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |
| 123 | (4-pyrimidin-5-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |
| 124 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-4-ylphenyl)methyl]urea |  |
| 125 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-3-ylphenyl)hydrazinecarboxamide |  |
| 126 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide |  |

Table 3

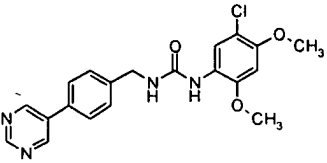
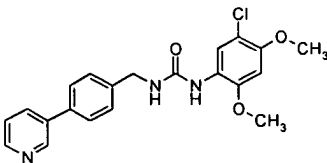
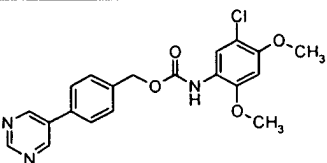
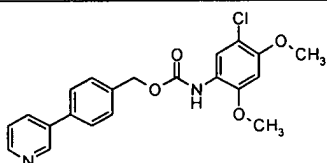
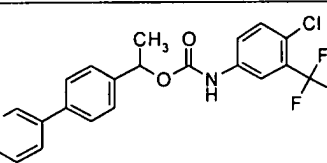
| Entry | Name | Structure |
|-------|---|--|
| 127 | N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyrimidin-5-ylphenyl)methyl]urea |  |
| 128 | N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea |  |
| 129 | (4-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate |  |
| 130 | (4-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate |  |
| 131 | 1-(4-pyridin-3-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |

Table 3

| Entry | Name | Structure |
|-------|---|-----------|
| 132 | 1-(4-pyrimidin-5-ylphenyl)ethyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 133 | N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(3-pyridin-3-ylphenyl)methyl]urea | |
| 134 | N-[5-chloro-2,4-bis(methoxy)phenyl]-N'-[(3-pyrimidin-5-ylphenyl)methyl]urea | |
| 135 | (3-pyridin-3-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate | |
| 136 | (3-pyrimidin-5-ylphenyl)methyl [5-chloro-2,4-bis(methoxy)phenyl]carbamate | |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 137 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-methyl-2-(3-pyrimidin-5-ylphenyl)hydrazinecarboxamide | |
| 138 | N-[4-fluoro-3-(trifluoromethyl)phenyl]-N'-[(4-pyridin-3-ylphenyl)methyl]urea | |
| 139 | N-{[3-(6-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea | |
| 140 | N-{[4-(6-aminopyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea | |
| 141 | N-{[3-(2-aminopyrimidin-5-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea | |

Table 3

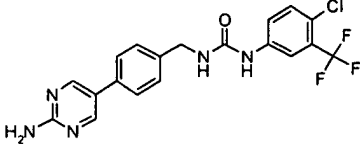
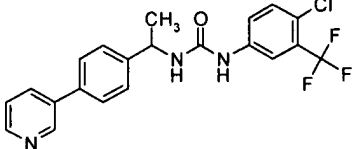
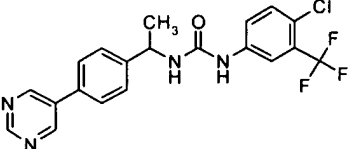
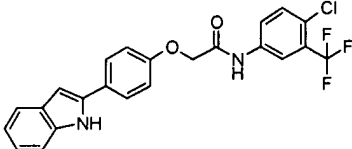
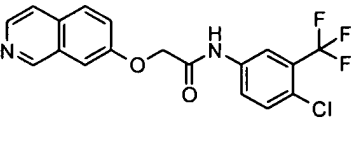
| Entry | Name | Structure |
|-------|---|--|
| 142 | N-{[4-(2-aminopyrimidin-5-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea |  |
| 143 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyridin-3-ylphenyl)ethyl]urea |  |
| 144 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[1-(4-pyrimidin-5-ylphenyl)ethyl]urea |  |
| 145 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[4-(1H-indol-2-yl)phenyl]oxy}acetamide |  |
| 146 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(isoquinolin-7-yloxy)acetamide |  |

Table 3

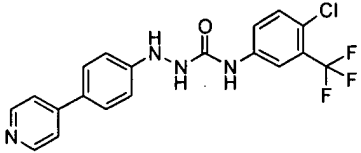
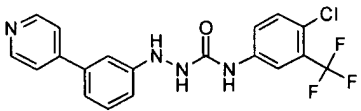
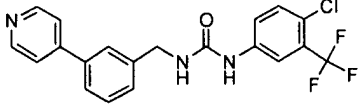
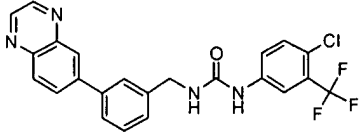
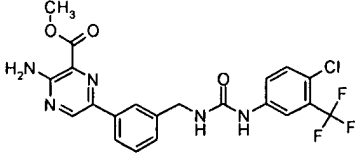
| Entry | Name | Structure |
|-------|--|--|
| 147 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(4-pyridin-4-ylphenyl)hydrazinecarboxamide |  |
| 148 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2-(3-pyridin-4-ylphenyl)hydrazinecarboxamide |  |
| 149 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-pyridin-4-ylphenyl)methyl]urea |  |
| 150 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-quinoxalin-6-ylphenyl)methyl]urea |  |
| 151 | methyl 3-amino-6-(3-{{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methyl}phenyl)pyrazine-2-carboxylate |  |

Table 3

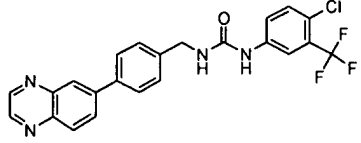
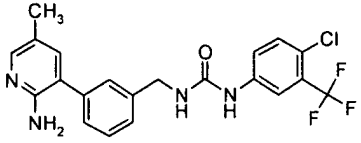
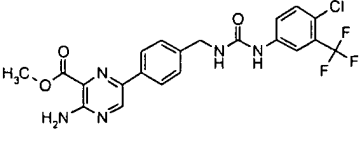
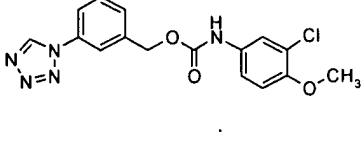
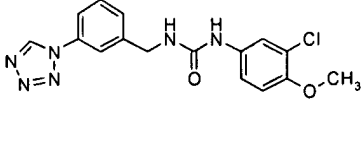
| Entry | Name | Structure |
|-------|---|--|
| 152 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-quinoxalin-6-ylphenyl)methyl]urea |  |
| 153 | N-{[3-(2-amino-5-methylpyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea |  |
| 154 | methyl 3-amino-6-(4-{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonylamino)methyl}phenyl)pyrazine-2-carboxylate |  |
| 155 | [3-(1H-tetrazol-1-yl)phenyl]methyl [3-chloro-4-(methyloxy)phenyl]carbamate |  |
| 156 | N-[3-chloro-4-(methyloxy)phenyl]-N'-{[3-(1H-tetrazol-1-yl)phenyl]methyl}urea |  |

Table 3

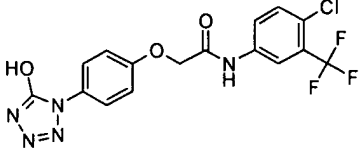
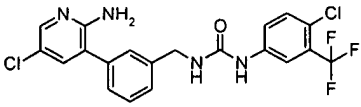
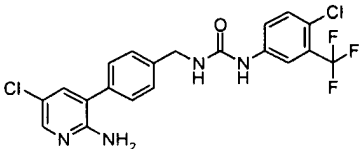
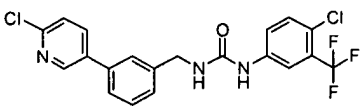
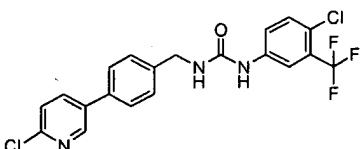
| Entry | Name | Structure |
|-------|--|--|
| 157 | N-[4-chloro-3-(trifluoromethyl)phenyl]-2- {[4-(5-hydroxy-1H-tetrazol-1- yl)phenyl]oxy}acetamide |  |
| 158 | N-{[3-(2-amino-5-chloropyridin-3- yl)phenyl]methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea |  |
| 159 | N-{[4-(2-amino-5-chloropyridin-3- yl)phenyl]methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea |  |
| 160 | N-{[3-(6-chloropyridin-3- yl)phenyl]methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea |  |
| 161 | N-{[4-(6-chloropyridin-3- yl)phenyl]methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea |  |

Table 3

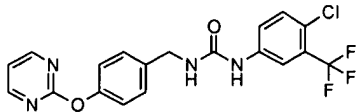
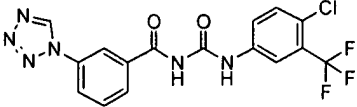
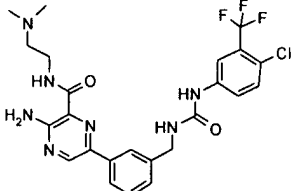
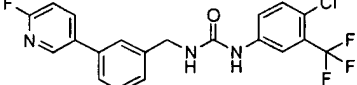
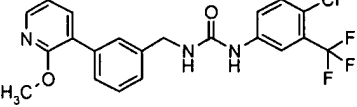
| Entry | Name | Structure |
|-------|---|--|
| 162 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(pyrimidin-2- yloxy)phenyl]methyl}urea |  |
| 163 | N-({[4-chloro-3- (trifluoromethyl)phenyl]amino}carbonyl)- 3-(1H-tetrazol-1-yl)benzamide |  |
| 164 | 3-amino-6-(3-{[({[4-chloro-3- (trifluoromethyl)phenyl]amino}carbonyl)a mino]methyl}phenyl)-N-[2- (dimethylamino)ethyl]pyrazine-2- carboxamide |  |
| 165 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[3-(6-fluoropyridin-3- yl)phenyl]methyl}urea |  |
| 166 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[2-(methyloxy)pyridin-3- yl]phenyl}methyl)urea |  |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 167 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(6-fluoropyridin-3- yl)phenyl]methyl}urea | |
| 168 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ([4-[2-(methyloxy)pyridin-3- yl]phenyl]methyl)urea | |
| 169 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(6-methylpyridin-3- yl)phenyl]methyl}urea | |
| 170 | N-{[4-(2-amino-5-fluoropyridin-3- yl)phenyl]methyl}-N'-[4-chloro-3- (trifluoromethyl)phenyl]urea | |
| 171 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[3-(6-methylpyridin-3- yl)phenyl]methyl}urea | |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 172 | N-{{4-(2-aminopyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea | |
| 173 | N-{{3-(2-aminopyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea | |
| 174 | [3-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 175 | [3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 176 | [3-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |

Table 3

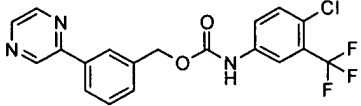
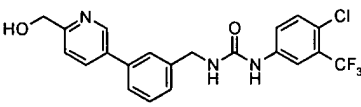
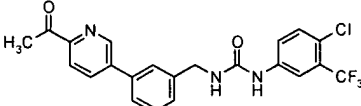
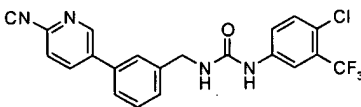
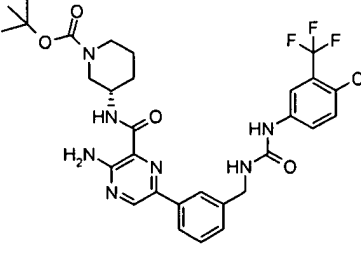
| Entry | Name | Structure |
|-------|---|--|
| 177 | (3-pyrazin-2-yl)phenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |
| 178 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[6-(hydroxymethyl)pyridin-3-yl]phenyl} methyl)urea |  |
| 179 | N-{{3-(6-acetylpyridin-3-yl)phenyl}methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea |  |
| 180 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{{3-(6-cyanopyridin-3-yl)phenyl}methyl} urea |  |
| 181 | 1,1-dimethylethyl (3S)-3-({[3-amino-6-(3-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazin-2-yl)carbonyl} amino)piperidine-1-carboxylate |  |

Table 3

| Entry | Name | Structure |
|-------|---|-----------|
| 182 | 3-amino-6-(3-{{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methyl}phenyl)-N-[(3S)-piperidin-3-yl]pyrazine-2-carboxamide | |
| 183 | 1,1-dimethylethyl (3S)-3-({[3-amino-6-(4-{{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methyl}phenyl]pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate | |
| 184 | 3-amino-6-(4-{{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methyl}phenyl)-N-[(3S)-piperidin-3-yl]pyrazine-2-carboxamide | |
| 185 | [3-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 186 | N-{[3-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea | |
| 187 | [6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 188 | [3-(1H-benzimidazol-2-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 189 | [3-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 190 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-({3-[5-(methylthio)pyridin-3-yl]phenyl}methyl)urea | |

Table 3

| Entry | Name | Structure |
|-------|---|-----------|
| 191 | [4-(6-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 192 | [4-(2-amino-5-fluoropyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 193 | [4-(2-aminopyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 194 | (4-pyrazin-2-ylphenyl)methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 195 | [4-(7H-pyrrolo[2,3-d]pyrimidin-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 196 | [4-(6-amino-2-methylpyridin-3-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 197 | [3-(1H-tetrazol-1-yl)phenyl]methyl 1,3-benzothiazol-2-ylcarbamate | |
| 198 | [3-(1H-tetrazol-1-yl)phenyl]methyl (5-bromopyridin-2-yl)carbamate | |
| 199 | (3-pyridin-3-ylphenyl)methyl (3,5-dimethylphenyl)carbamate | |
| 200 | (3-pyridin-3-ylphenyl)methyl [5-chloro-2-(methyloxy)phenyl]carbamate | |

Table 3

| Entry | Name | Structure |
|-------|---|-----------|
| 201 | [4-(1H-tetrazol-1-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate | |
| 202 | (3-pyrimidin-5-ylphenyl)methyl [5-chloro-2-(methoxy)phenyl]carbamate | |
| 203 | (4-pyrimidin-5-ylphenyl)methyl (3,4-dimethylphenyl)carbamate | |
| 204 | (3-pyridin-3-ylphenyl)methyl (3,4-dimethylphenyl)carbamate | |
| 205 | 1,1-dimethylethyl 3-({[3-amino-6-(3-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl)amino]methyl}phenyl)pyrazin-2-yl)carbonyl}amino)piperidine-1-carboxylate | |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 206 | 1,1-dimethylethyl 3-({[3-amino-6-(4- {{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methyl}phenyl)pyrazin-2-yl]carbonyl}amino)piperidine-1-carboxylate | |
| 207 | 3-amino-6-(3-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methyl}phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide | |
| 208 | 3-amino-6-(4-({[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methyl}phenyl)-N-piperidin-3-ylpyrazine-2-carboxamide | |
| 209 | 1,1-dimethylethyl 4-({[3-amino-6-(3- {{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methyl}phenyl)pyrazin-2-yl]carbonyl}piperazine-1-carboxylate | |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 210 | 1,1-dimethylethyl 4-{{[3-amino-6-(4- {{{[4-chloro-3-(trifluoromethyl)phenyl]amino}carbonyl]amino}methyl]phenyl}pyrazin-2-yl]carbonyl}piperazine-1-carboxylate | |
| 211 | N-({3-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea | |
| 212 | N-({4-[5-amino-6-(piperazin-1-ylcarbonyl)pyrazin-2-yl]phenyl}methyl)-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea | |
| 213 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{{[3-(1H-pyrazol-4-yl)phenyl]methyl}urea | |

Table 3

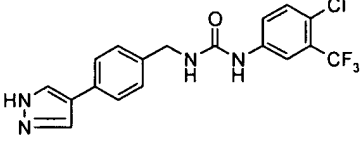
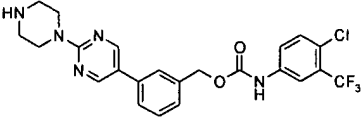
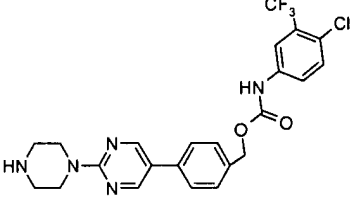
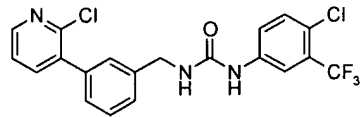
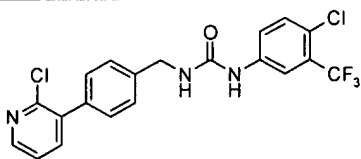
| Entry | Name | Structure |
|-------|--|--|
| 214 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-{[4-(1H-pyrazol-4-yl)phenyl]methyl}urea |  |
| 215 | [3-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |
| 216 | [4-(2-piperazin-1-ylpyrimidin-5-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |
| 217 | N-{[3-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea |  |
| 218 | N-{[4-(2-chloropyridin-3-yl)phenyl]methyl}-N'-[4-chloro-3-(trifluoromethyl)phenyl]urea |  |

Table 3

| Entry | Name | Structure |
|-------|--|-----------|
| 219 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[3-(2-fluoropyridin-3-yl)phenyl]methyl}urea | |
| 220 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- {[4-(2-fluoropyridin-3-yl)phenyl]methyl}urea | |
| 221 | [3-(1H-tetrazol-1-yl)phenyl]methyl [3-(trifluoromethyl)phenyl]carbamate | |
| 222 | [3-(1H-tetrazol-1-yl)phenyl]methyl [6-(trifluoromethyl)pyridin-2-yl]carbamate | |
| 223 | [3-(1H-tetrazol-1-yl)phenyl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate | |

Table 3

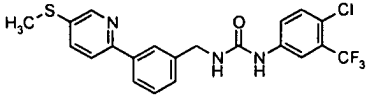
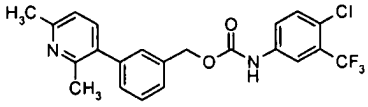
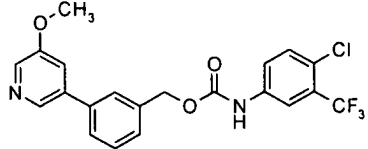
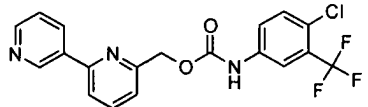
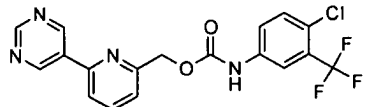
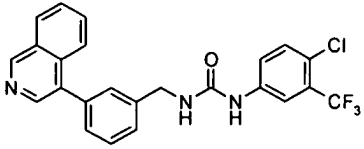
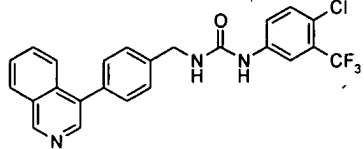
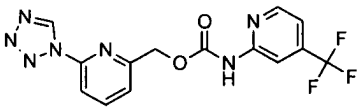
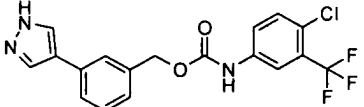
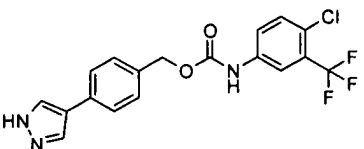
| Entry | Name | Structure |
|-------|--|--|
| 224 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'- ({3-[5-(methylthio)pyridin-2- yl]phenyl}methyl)urea |  |
| 225 | [3-(2,6-dimethylpyridin-3- yl)phenyl]methyl [4-chloro-3- (trifluoromethyl)phenyl]carbamate |  |
| 226 | {3-[5-(methyloxy)pyridin-3- yl]phenyl}methyl [4-chloro-3- (trifluoromethyl)phenyl]carbamate |  |
| 227 | 2,3'-bipyridin-6-ylmethyl [4-chloro-3- (trifluoromethyl)phenyl]carbamate |  |
| 228 | (6-pyrimidin-5-ylpyridin-2-yl)methyl [4- chloro-3- (trifluoromethyl)phenyl]carbamate |  |

Table 3

| Entry | Name | Structure |
|-------|---|--|
| 229 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(3-isoquinolin-4-ylphenyl)methyl]urea |  |
| 230 | N-[4-chloro-3-(trifluoromethyl)phenyl]-N'-[(4-isoquinolin-4-ylphenyl)methyl]urea |  |
| 231 | [6-(1H-tetrazol-1-yl)pyridin-2-yl]methyl [4-(trifluoromethyl)pyridin-2-yl]carbamate |  |
| 232 | [3-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |
| 233 | [4-(1H-pyrazol-4-yl)phenyl]methyl [4-chloro-3-(trifluoromethyl)phenyl]carbamate |  |

28. (previously presented) A pharmaceutical composition comprising the compound according to claim 1 and a pharmaceutically acceptable carrier.

29. (cancelled)

30. (withdrawn) A method for modulating the *in-vivo* activity of a kinase, the method comprising administering to a subject an effective amount of the compound according to claim 1 or a compound selected from N-naphthalen-1-yl-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[4-(phenoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(3,4-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,3-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,4-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,5-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(3,5-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,6-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,6-diethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2-(methoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2-(ethoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[3-(ethoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2,4-bis(methoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2,3-dichlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-bromophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(2-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-N-[2-(trifluoromethyl)phenyl] acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl}oxy}-N-[3-(trifluoromethyl)phenyl] acetamide, methyl 4-[[{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetyl]amino] benzoate, ethyl 4-[[{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetyl]amino] benzoate, 3-[[{{3-(1H-tetrazol-1-yl)phenyl}oxy}acetyl]amino] benzoic acid, N-[3-(methoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[4-(methoxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-chlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl}oxy} acetamide, N-(4-aminophenyl)-2-{{3-(1H-

tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide.

31. (withdrawn) The method according to claim 30, wherein the kinase is c-Kit.

32. (withdrawn) The method according to claim 31, wherein modulating the *in vivo* activity of c-Kit comprises inhibition of c-Kit.

33. (withdrawn) A method of treating diseases or disorders associated with uncontrolled, abnormal, and/or unwanted cellular activities, the method comprising administering, to a mammal in need thereof, a therapeutically effective amount of the compound or the pharmaceutical composition as described in claim 1 or a compound, or a pharmaceutical composition comprising said compound, selected from N-naphthalen-1-yl-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethyl-phenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl) acetamide, N-(2-ethyl-phenyl)-2-{{3-(1H-tetrazol-1-yl) phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(methyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyl-oxy)phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)-phenyl]-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-bromophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{{3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{{3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-[[{{3-(1H-tetrazol-1-

yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-([3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-([3-(1H-tetrazol-1-yl)phenyl]oxy} acetyl) amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl] oxy} acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide.

34. (withdrawn) A method of screening for modulators of c-Kit, the method comprising combining the compound according to claim 1 or a compound selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl] -2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethyl-phenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl] oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl) phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy) phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-

N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl amino benzoate, ethyl 4-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl amino benzoate, 3-([3-(1H-tetrazol-1-yl)phenyl]oxy)acetyl amino benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and at least one candidate agent and determining the effect of the candidate agent on c-Kit activity.

35. (withdrawn) A method of inhibiting proliferative activity in a cell, the method comprising administering an effective amount of a composition comprising the compound according to claim 1 or a compound selected from N-naphthalen-1-yl-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(phenyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,4-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(3,5-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-dimethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-(2,4,6-trimethylphenyl)acetamide, N-(2-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-ethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,6-diethylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[3-(ethyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2,4-bis(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(dimethylamino)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2,3-dichlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-chloro-3-methylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(4-bromophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-(2-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide,

acetamide, N-(4-fluorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[2-(trifluoro-methyl)phenyl] acetamide, 2-{[3-(1H-tetrazol-1-yl)phenyl]oxy}-N-[3-(trifluoromethyl) phenyl] acetamide, methyl 4-[(3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, ethyl 4-[(3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoate, 3-[(3-(1H-tetrazol-1-yl)phenyl]oxy}acetyl)amino] benzoic acid, N-[3-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-(methyloxy)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[2-chloro-5-(trifluoromethyl)phenyl]-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, N-[4-chloro-3-(trifluoromethyl)phenyl]-2-{[3-(4H-1,2,4-triazol-4-yl)phenyl]oxy} acetamide, N-(4-chlorophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl] oxy} acetamide, N-(4-aminophenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, and N-(4-acetylphenyl)-2-{[3-(1H-tetrazol-1-yl)phenyl]oxy} acetamide, to a cell or a plurality of cells.